

Improving the AMPT Model for Isobar Collisions

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National
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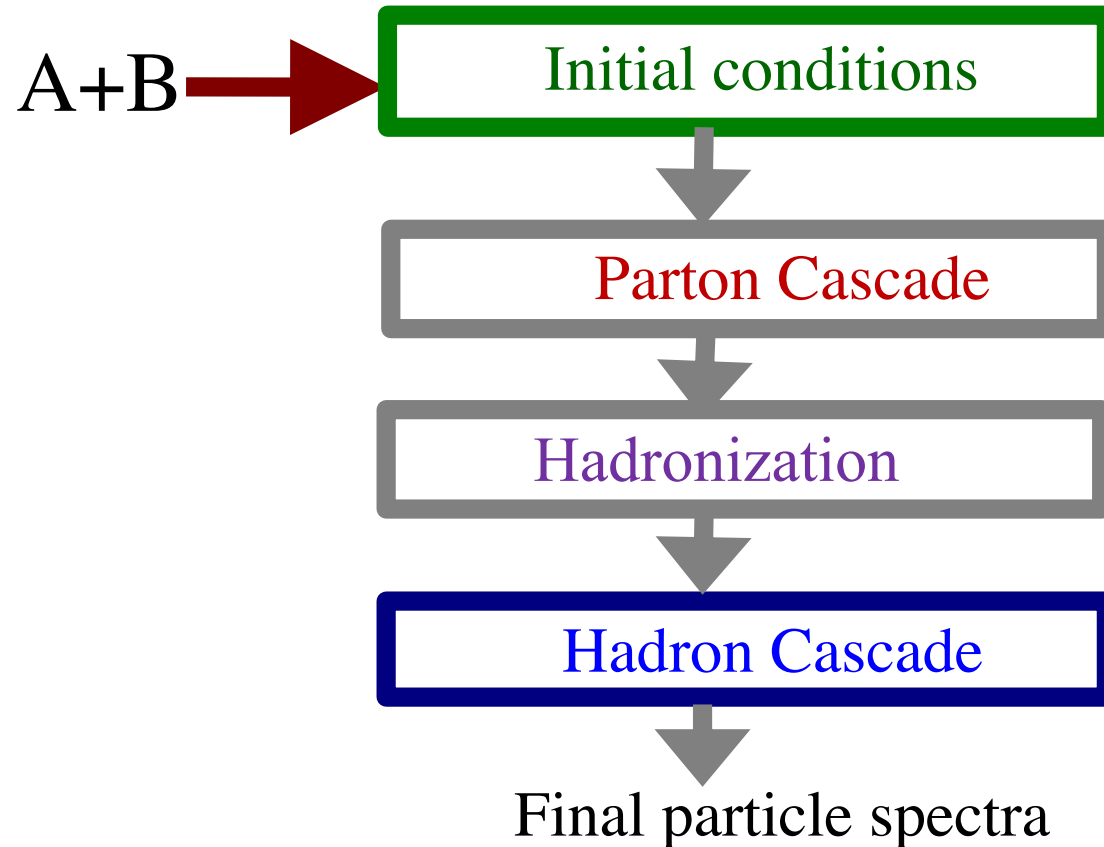
Outline

- Introduction of the AMPT model
 - Improvements relevant for isobar collisions
 - Incorporation of nucleus of different shapes
 - Local nuclear scaling and self-consistent system size dependence
 - The new quark coalescence model
 - Implementation of charge conservation
 - Further developments
 - Summary
- Partly based on a mini-review of recent AMPT developments on Nucl. Sci. Tech. (2021) and collaborations with Guo-Liang Ma, Bedangadas Mohanty, Md. Rihan Haque, Fuqiang Wang, Chao Zhang, Liang Zheng, et al.

A Multi-Phase Transport (AMPT)

Constructed as a self-contained kinetic description of heavy ion collisions

- evolves the system from fluctuating initial conditions to final observables;
- produces particles of all flavours at all P_T & y ;
- includes non-equilibrium initial condition & dynamics.



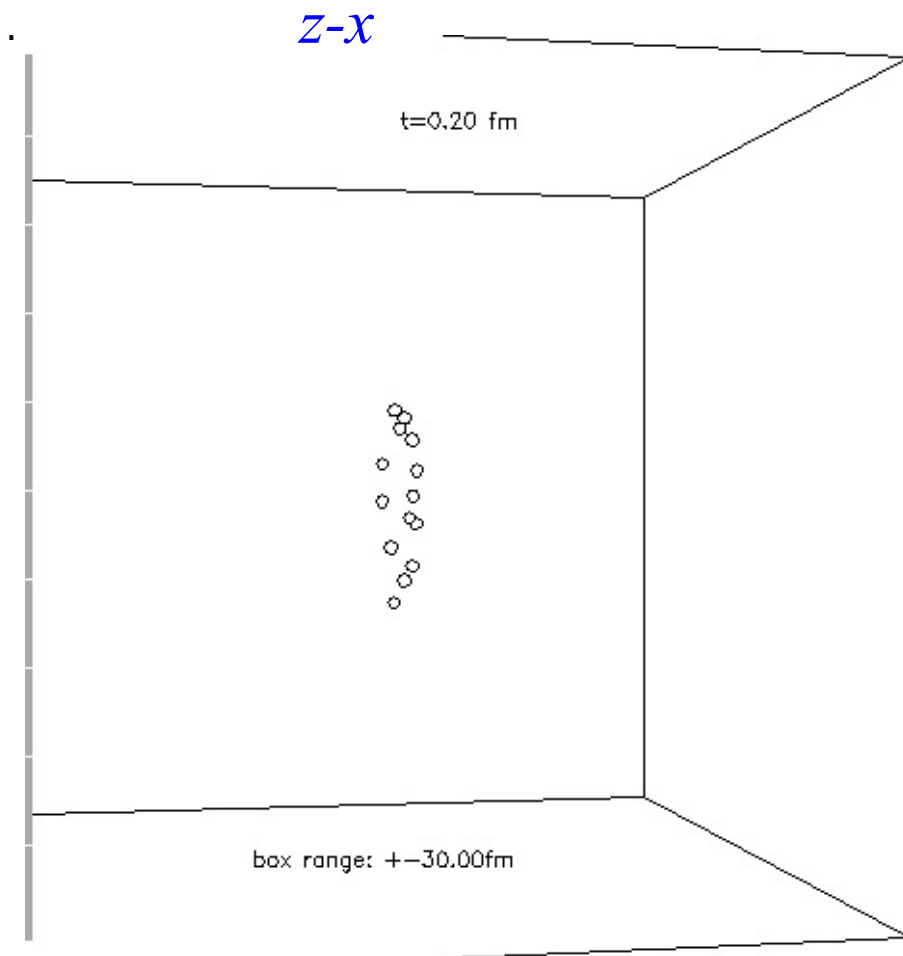
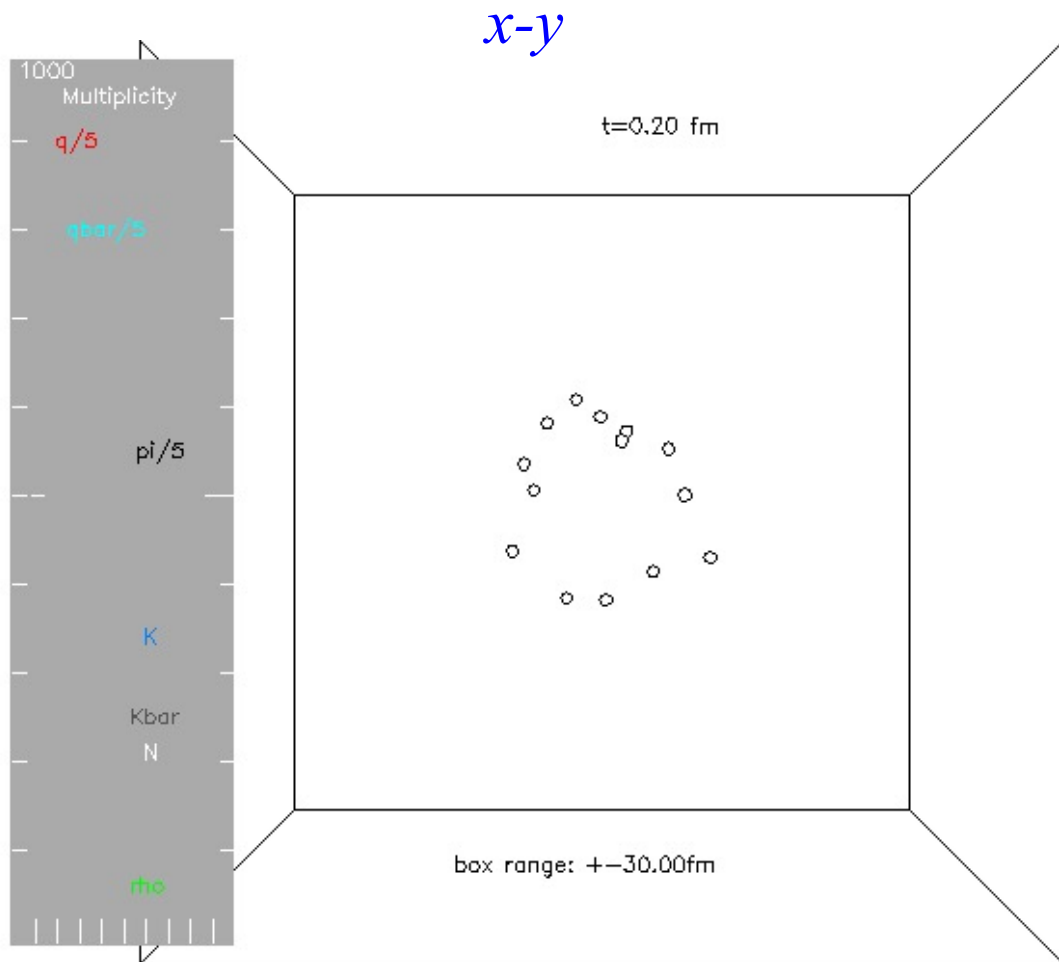
Source codes online since 2004,
available at the ECU website

<http://myweb.ecu.edu/linz/ampt/>

ZWL, Ko, Li, Zhang & Pal, PRC (2005);
ZWL & Zheng, NST (2021).

A Multi-Phase Transport (AMPT)

Time evolution of 1 central Au+Au event at 200A GeV
from AMPT-SM (*the String Melting version, applicable at high energies*):



Animations at the ECU website
<http://myweb.ecu.edu/linz/ampt/>

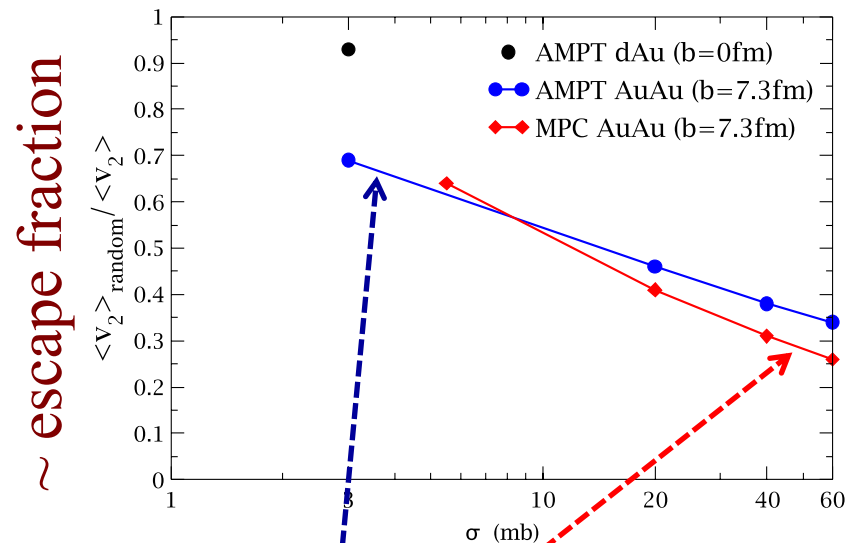
Transport models for finite systems

- **For large systems at very high energies:**
transport model approaches **hydrodynamics**,
transport model (particles, scatterings, microscopic picture) are complementary to **hydrodynamics-based models** ($T_{\mu\nu}$, EoS, transport coefficients).
- **For finite/small systems at finite energies:**
non-equilibrium effects can be important,
important to develop **transport model/kinetic theory** & compare with **hydrodynamics** to understand physics including collectivity of finite size systems.

Heiselberg & Levy, PRC (1999), Borghini et al. EPJC (2018), Kurkela et al. PLB (2018) & EPJC (2019), ...

On example is
the escape mechanism:
interaction-induced response
to anisotropic geometry
from kinetic theory.

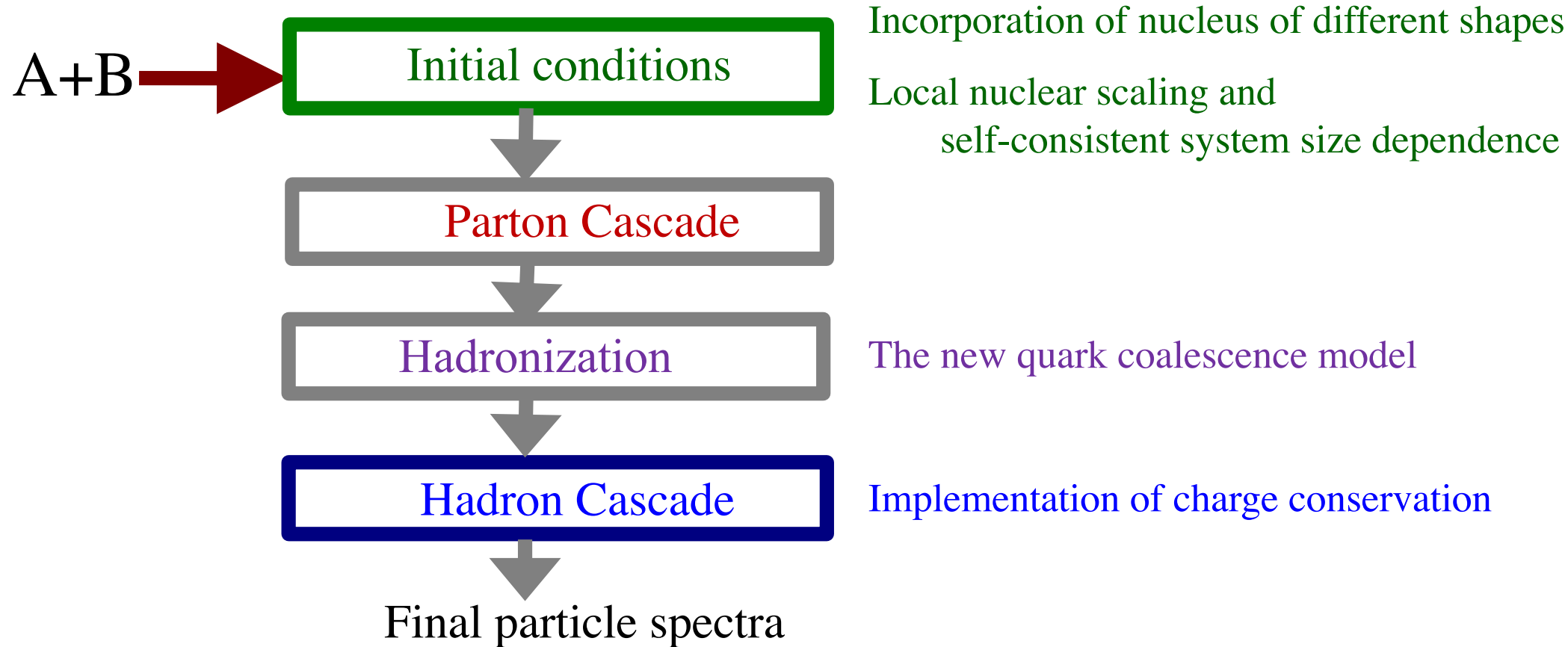
L He et al. PLB (2016);
ZWL et al. NPA (2016);
H.L. Li et al. PRC (2019).



- At very large opacity (large system/energy/ σ), **hydrodynamic collective flow** will dominate v_2
- **Escape mechanism** dominates v_2 for small systems & even semi-central AuAu @200 GeV.

Improvements relevant for isobar collisions

They cover the following AMPT components:



ZWL & Zheng, NST (2021)

Improvements not covered in this talk: updates with modern $nPDFs$, heavy flavor productions, finite nuclear thickness, pythia8 initial condition with sub-nucleon structure, parton cascade.

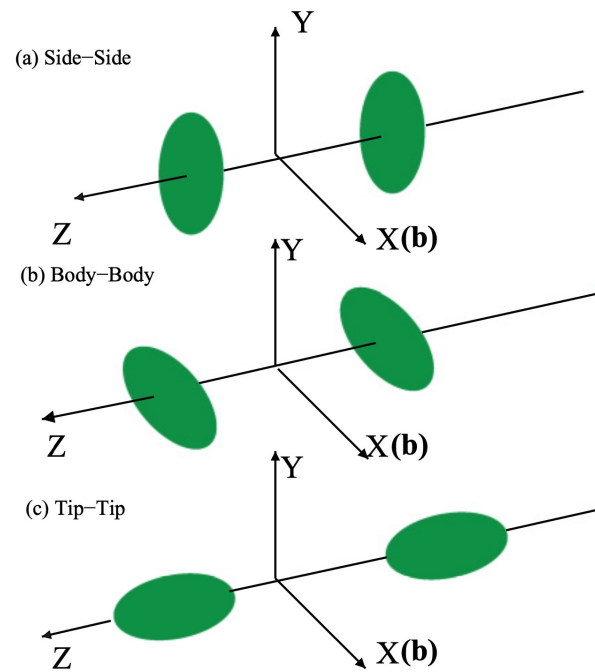
Incorporation of nucleus of different shapes

Rihan Haque, ZWL & Mohanty, PRC (2012)

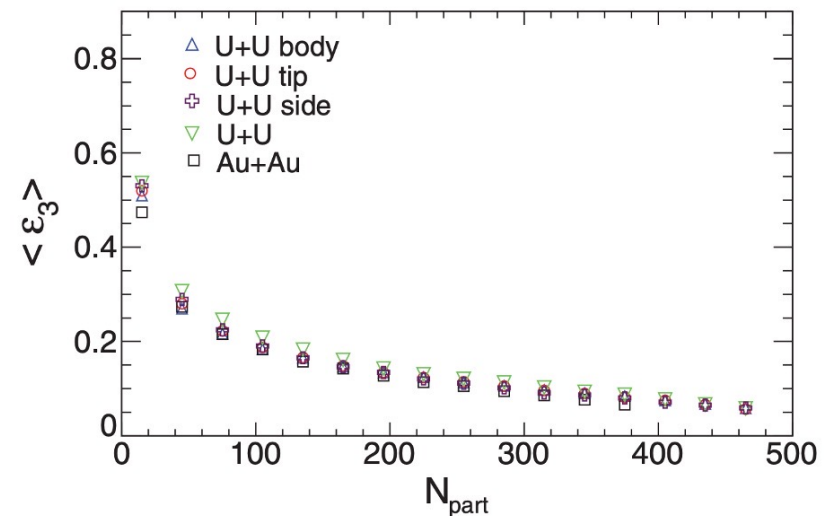
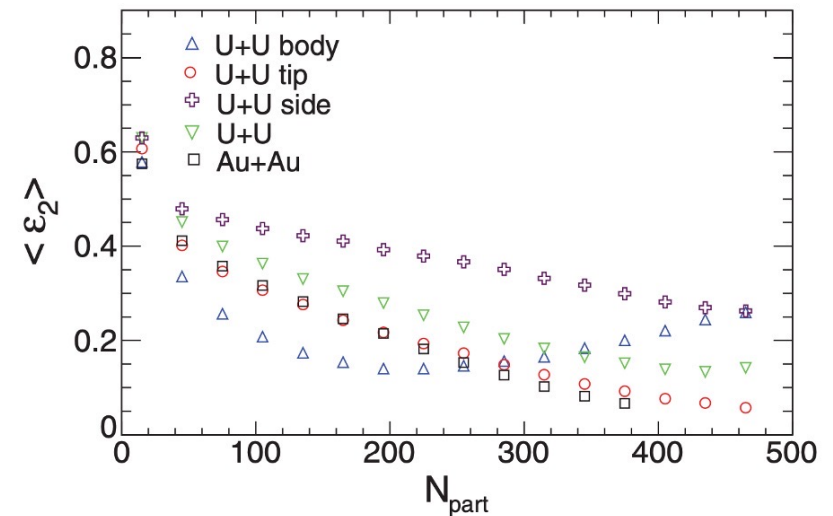
Nucleon density distribution of ^{238}U in AMPT is parametrized as deformed Woods-Saxon:

$$\rho = \frac{\rho_0}{1 + \exp([r - R']/a)},$$

$$R' = R[1 + \beta_2 Y_2^0(\theta) + \beta_4 Y_4^0(\theta)]$$

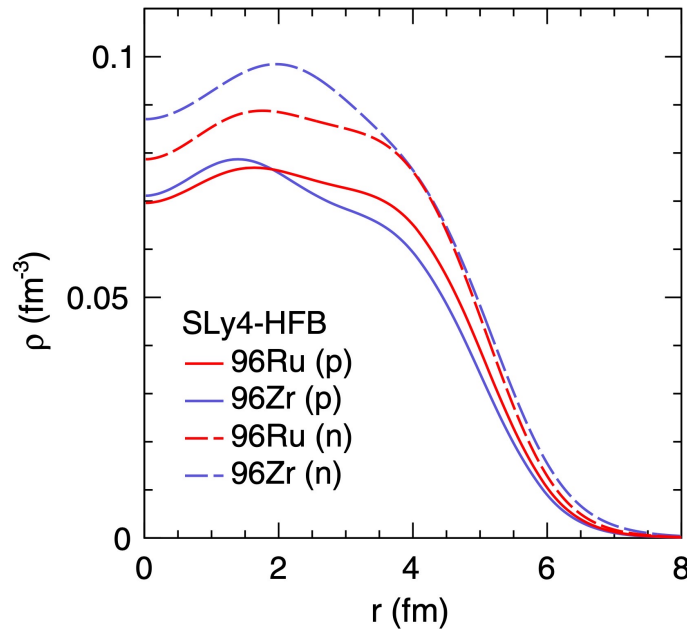


Eccentricity ε_2 and triangularity ε_3 :
 ε_2 and v_2 are very sensitive to orientation.



Incorporation of nucleus of different shapes

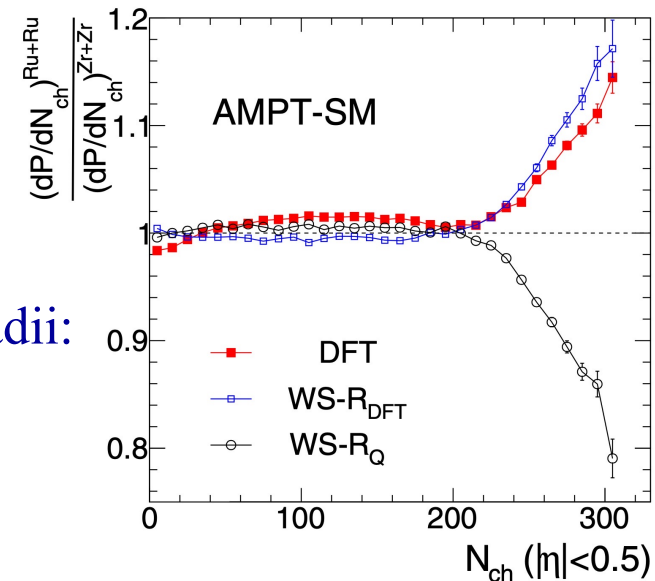
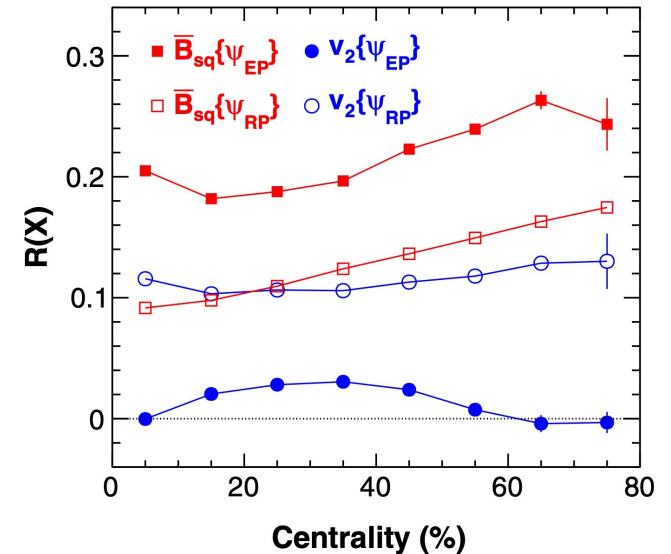
For isobar nuclei in the AMPT initial condition, protons & neutrons are sampled according to radial density distributions from density functional theory (DFT):



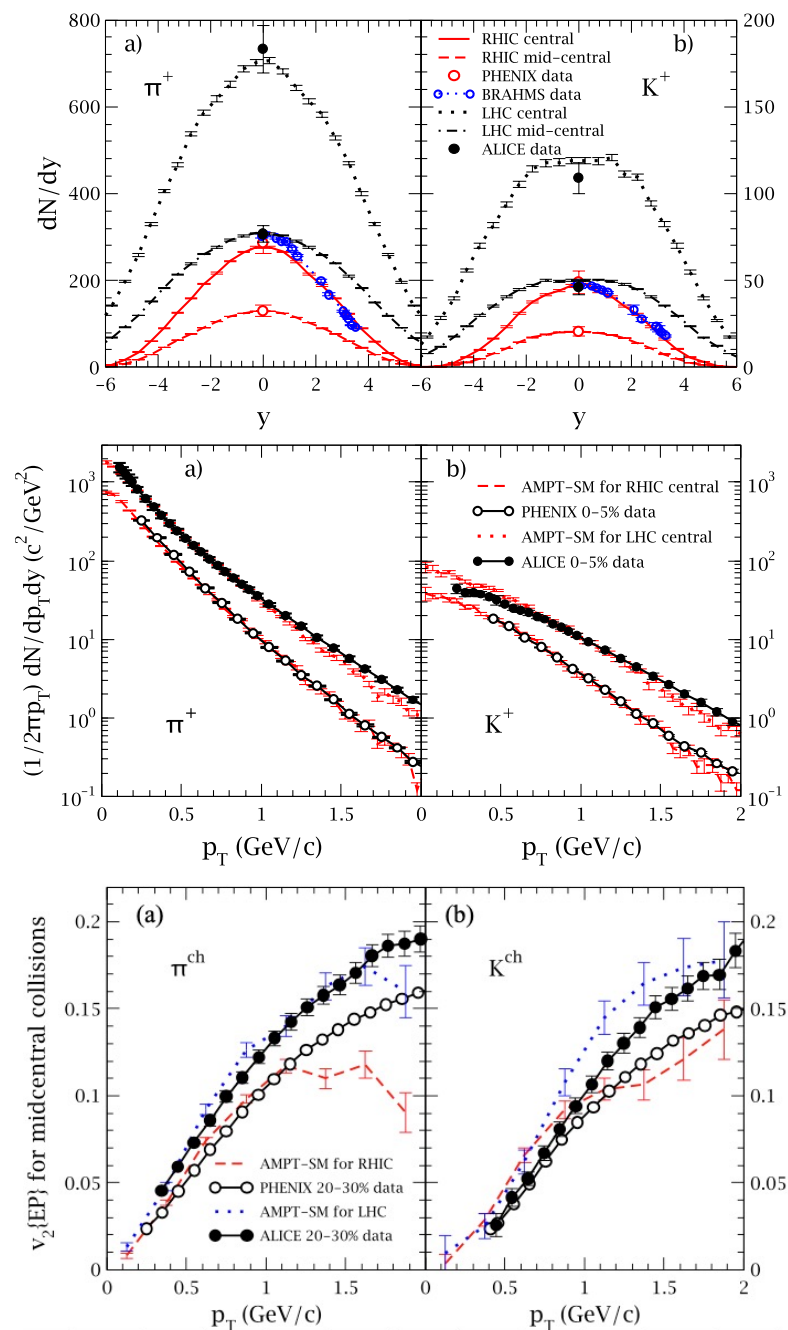
H.J. Xu et al. PRL (2018);
H.J. Xu's talk on Tuesday

Ratio of N_{ch} distributions
probes the nuclear mass radii:
H.L. Li et al. PRC (2018)

Sizable v_2 difference is found for isobar systems:



Local nuclear scaling and self-consistent size dependence



String Melting AMPT describes flows and HBT
but used to (*before 2014*) fail badly in hadron spectra.

We later realized that the model
can simultaneously describe dN/dy , p_T -spectra & v_2
(for bulk matter at low- p_T) in central and
mid-central high energy AA collisions,
as long as a very small Lund b_L parameter is used.

ZWL, PRC (2014)

Lund symmetric string fragmentation function:

$$f(z) \propto z^{-1} (1-z)^{a_L} e^{-b_L m_T^2/z}$$

b_L typical values (in $1/\text{GeV}^2$):
0.5 (AMPT-def), ~ 0.58 (PYTHIA6.2), 0.9 (HIJING1)

**$b_L \sim 0.15$ is needed for AMPT-SM for high energy AA,
corresponds to a much higher string tension:**

$$\kappa \propto \frac{1}{b_L(2+a_L)}$$

ZWL et al. PRC (2005)

Local nuclear scaling and self-consistent size dependence

Different values of b_L are needed for pp and central AA, same for the minijet cutoff scale p_0 (related to saturation scale Q_s in the AMPT model updated with modern $nPDFs$).

C Zhang et al. PRC (2019);
Zheng et al. PRC (2020)

We propose to scale them with local nuclear densities:

C Zhang et al.
PRC (2021)

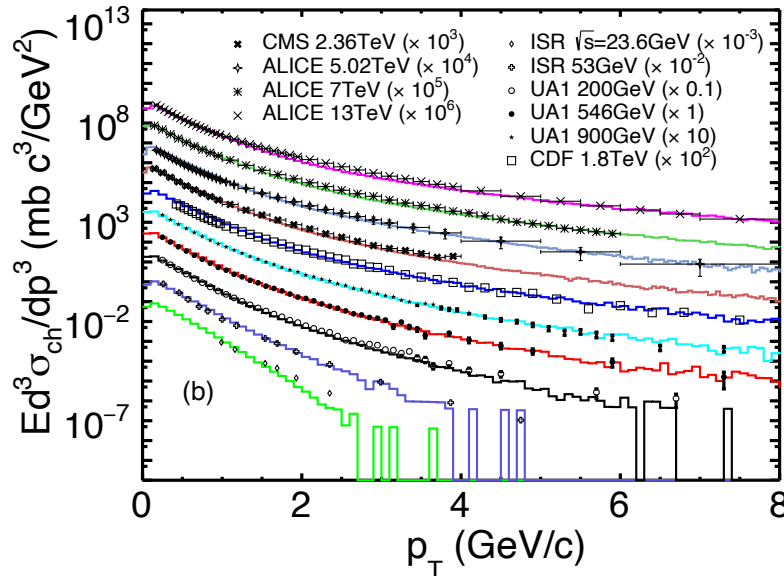
$$b_L(s_A, s_B, s) = \frac{b_L^{pp}}{[\sqrt{T_A(s_A)T_B(s_B)}/T_p]^{\beta(s)}}$$

$$p_0(s_A, s_B, s) = p_0^{pp}(s)[\sqrt{T_A(s_A)T_B(s_B)}/T_p]^{\alpha(s)}$$

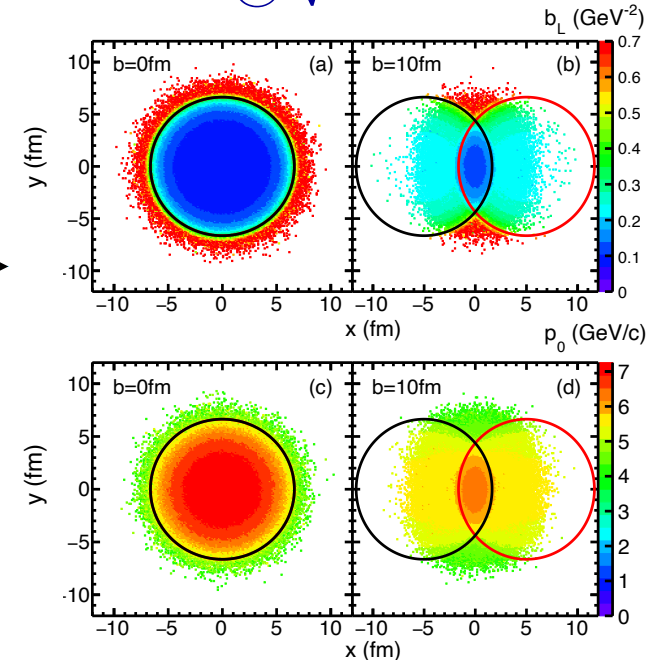
Similar geometric form preferred for initial entropy deposition:

Bernhard et al. PRC (2016)

We fit charged hadrons in pp to determine $b_L^{pp}=0.7$, then used central AuAu/PbPb data to get $\alpha(s)$, $\beta(s)$:



Pb+Pb @ $\sqrt{s}=5.02$ ATeV

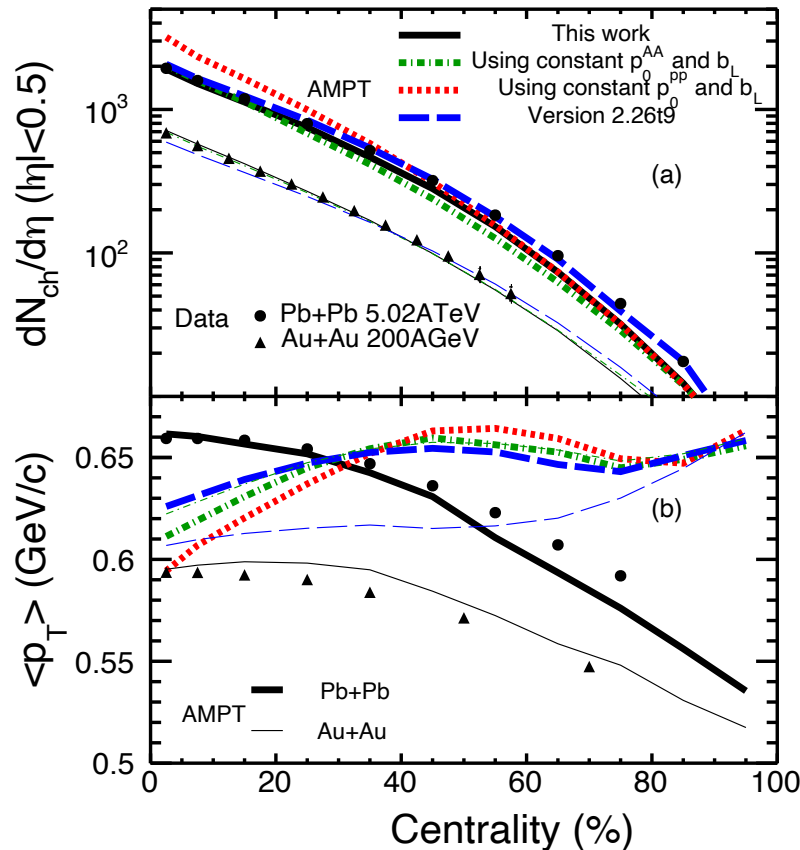


Local nuclear scaling and self-consistent size dependence

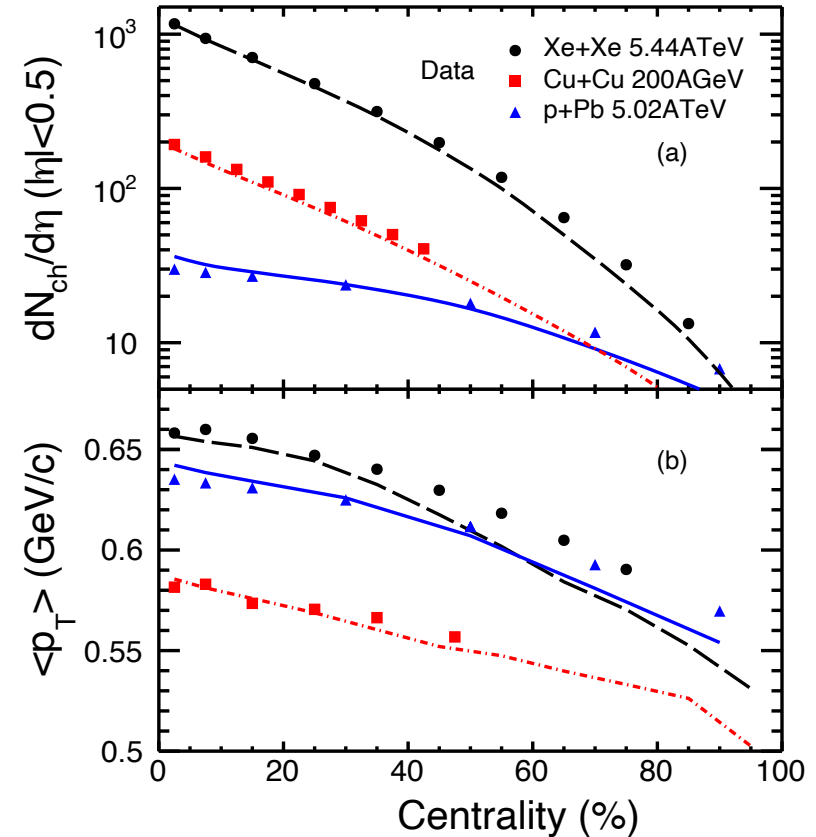
The scaling allows AMPT-SM to self-consistently describe the system size dependence, including centrality dependences of AuAu & PbPb and smaller systems.

C Zhang et al. PRC (2021)

Centrality dependences of $\langle p_T \rangle$ are now reasonable, much better than public AMPT (v2.26t9)



Also works for smaller systems:



Key input parameters of AMPT: a_L b_L p_0

No longer free parameters

σ (parton cross section)

$\rightarrow \eta \dots$ can be better studied

The new quark coalescence model

Old Coalescence

(in public AMPT)

→
 $q_{\bar{B}} \quad q_m \quad q_B \quad \bar{q}_m \dots\dots$

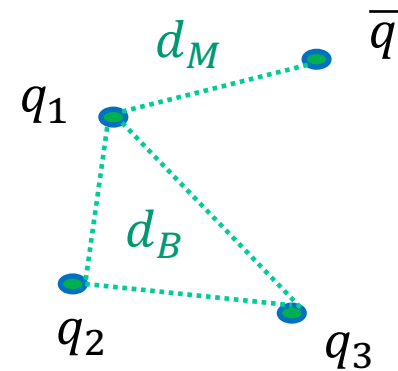
There is an artificial constraint that forces the separate conservation of the numbers of mesons, baryons and antibaryons for each event (*while only net-baryon conservation is necessary*):

- Quarks from the melting of mesons search all antiquarks and choose the closest antiquark to form mesons.
- Then quarks from the melting of baryons search all remaining quarks and choose the closest two quarks to form baryons (*same for anti-baryons*).

New coalescence

Y. He & ZWL, PRC (2017)

We remove the artificial constraint, **quarks now have freedom to form either meson or baryon:**



For example, for a quark q_1 :

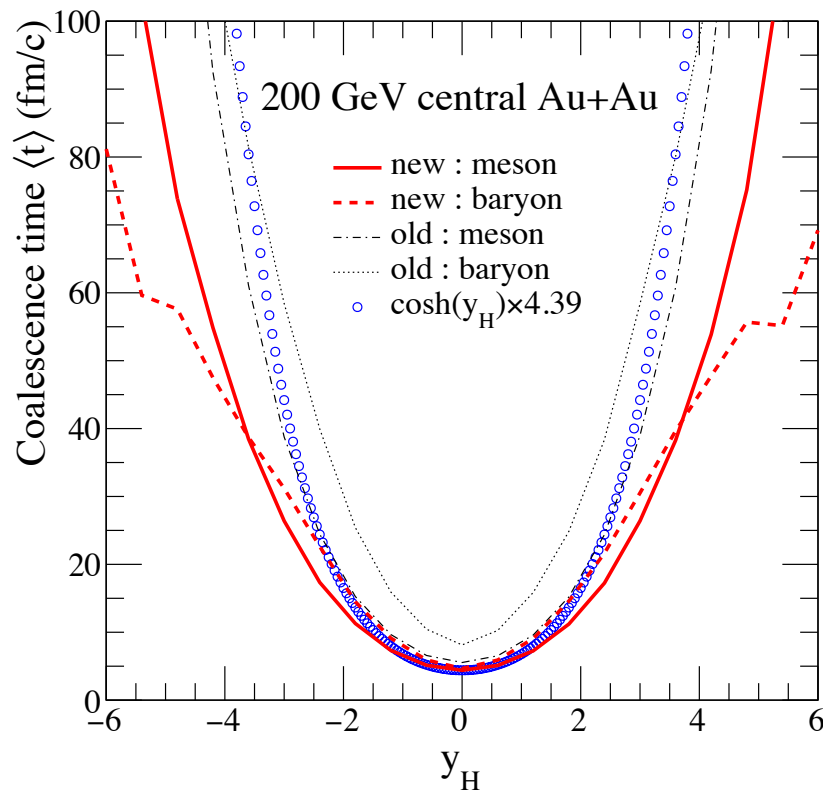
d_M : closest distance to an antiquark (*in rest frame*)

d_B : average distance among the 3 quarks
after finding closest q_2 & q_3

If $d_B < d_M * r_{BM}$, q_1 will coalesce to a baryon;
otherwise, q_1 will coalesce to a meson.

→ Single coalescence parameter r_{BM} , ~ 0.6

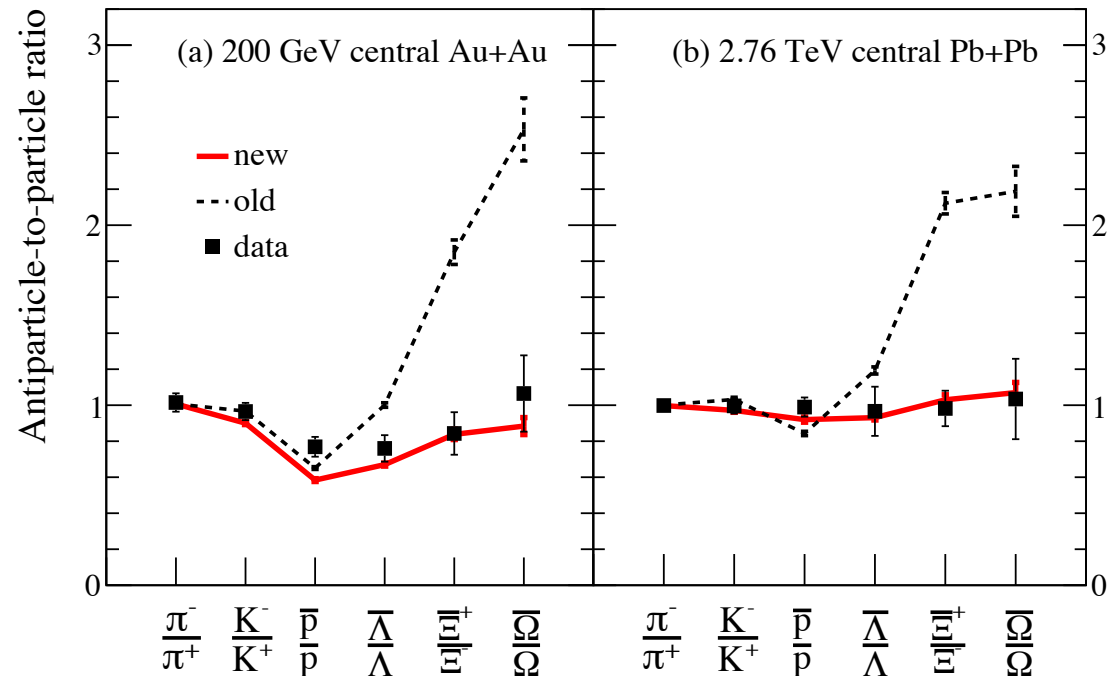
The new quark coalescence model



Y. He & ZWL, PRC (2017)

New quark coalescence is more efficient, especially for (anti)baryons.

Improves (anti)baryon observables, including p & pbar yield & p_T -spectra and multi-strange Bbar/B ratios:



Implementation of charge conservation

Why is charge conservation violated in public AMPT?

First reason: The hadron cascade has K^+ and K^- as explicit particles, but not K^0 and $K^0\text{-bar}$.

Currently, to let all kaons interact:

- before hadron cascade, we **change** K^0 to K^+ (also: $K^0\text{-bar}$ to K^-)
- after hadron cascade, we **change** half of final K^+ into K^0 .

Second reason:

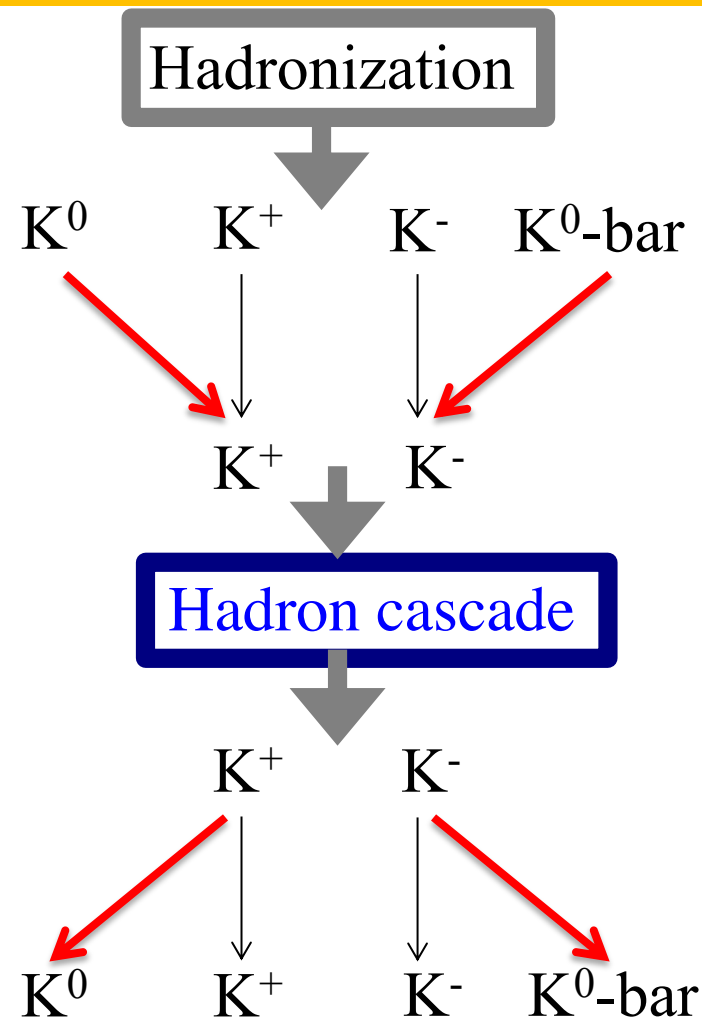
Many reactions in the hadron cascade are not implemented for each possible isospin/charge state

→ **Need to identify & correct each such reaction while respecting detailed balance relations**

This work has been done in a version of AMPT:

Z.W. Lin & G.L. Ma, unpublished (2018).

This charge-conserved version has been shared with some colleagues for CME studies: Tang, CPC (2020); Choudhury et al. EPJC (2020).



Further developments relevant for isobar collisions

- Further improve the initial condition from HIJING:

There is a small p_z -asymmetry of hadrons in symmetric collisions:
this leads to rapidity-asymmetry of final state hadrons,
especially antibaryons from public AMPT-SM at low energies
(*thanks to H.Z. Huan's group for pointing out this issue*).

There is an artificial ordering of initial protons and neutrons
in the nucleus along the z direction when sampling Woods-Saxon
(*protons take \sim half of the z range & neutrons take the other \sim half*):
this affects the initial net-charge distribution
and could affect charge-dependent observables.

- Self-consistent parton transport under electromagnetic fields
- Update the public AMPT model with recent improvements

Summary

AMPT provides a self-contained kinetic description of heavy ion collisions

Recent improvements make AMPT more versatile and reliable:

- **Incorporation of deformed or arbitrary nucleon distributions**
enables studies of isobar collisions.
- **New quark coalescence**
improves baryon & antibaryon productions.
- **Implementation of charge conservation**
improves studies of charge-dependent observables.
- **Local nuclear scaling of initial condition parameters**
significantly reduces uncertainty from free model parameters
and enables us to focus on QGP properties like
 σ (parton cross section, or m_D) and its T-dependence $\leftrightarrow \eta(T)$

Thanks for your attention!